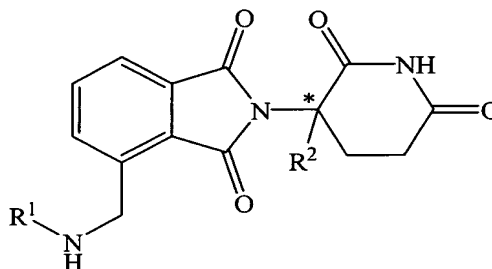


Amendments to the Claims

The following listing of claims will replace all prior versions, and listings, of claims in the application:

1-9. (Canceled).

10. (Previously presented) A compound having the formula:



wherein:

R^1 is H, $(\text{C}_1\text{--C}_8)\text{alkyl}$, $(\text{C}_3\text{--C}_7)\text{cycloalkyl}$, $(\text{C}_2\text{--C}_8)\text{alkenyl}$, $(\text{C}_2\text{--C}_8)\text{alkynyl}$, benzyl, aryl, $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_1\text{--C}_6)\text{heterocycloalkyl}$, $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_2\text{--C}_5)\text{heteroaryl}$, $\text{C}(\text{O})\text{R}^3$, $\text{C}(\text{S})\text{R}^3$, $\text{C}(\text{O})\text{OR}^4$, $(\text{C}_1\text{--C}_8)\text{alkyl--N}(\text{R}^6)_2$, $(\text{C}_1\text{--C}_8)\text{alkyl--OR}^5$, $(\text{C}_1\text{--C}_8)\text{alkyl--C}(\text{O})\text{OR}^5$, $\text{C}(\text{O})\text{NHR}^3$, $\text{C}(\text{S})\text{NHR}^3$, $\text{C}(\text{O})\text{NR}^3\text{R}^{3'}$, $\text{C}(\text{S})\text{NR}^3\text{R}^{3'}$ or $(\text{C}_1\text{--C}_8)\text{alkyl--O}(\text{CO})\text{R}^5$;

R^2 is H or $(\text{C}_1\text{--C}_8)\text{alkyl}$;

R^3 and $\text{R}^{3'}$ are independently $(\text{C}_1\text{--C}_8)\text{alkyl}$, $(\text{C}_3\text{--C}_7)\text{cycloalkyl}$, $(\text{C}_2\text{--C}_8)\text{alkenyl}$, $(\text{C}_2\text{--C}_8)\text{alkynyl}$, benzyl, aryl, $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_1\text{--C}_6)\text{heterocycloalkyl}$, $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_2\text{--C}_5)\text{heteroaryl}$, $(\text{C}_0\text{--C}_8)\text{alkyl--N}(\text{R}^6)_2$, $(\text{C}_1\text{--C}_8)\text{alkyl--OR}^5$, $(\text{C}_1\text{--C}_8)\text{alkyl--C}(\text{O})\text{OR}^5$, $(\text{C}_1\text{--C}_8)\text{alkyl--O}(\text{CO})\text{R}^5$, or $\text{C}(\text{O})\text{OR}^5$;

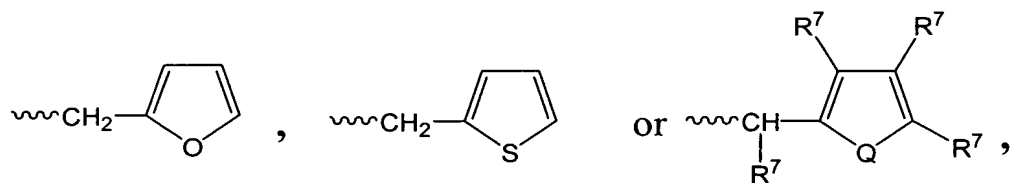
R^4 is $(\text{C}_1\text{--C}_8)\text{alkyl}$, $(\text{C}_2\text{--C}_8)\text{alkenyl}$, $(\text{C}_2\text{--C}_8)\text{alkynyl}$, $(\text{C}_1\text{--C}_4)\text{alkyl--OR}^5$, benzyl, aryl, $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_1\text{--C}_6)\text{heterocycloalkyl}$, or $(\text{C}_0\text{--C}_4)\text{alkyl--}(\text{C}_2\text{--C}_5)\text{heteroaryl}$;

R^5 is $(\text{C}_1\text{--C}_8)\text{alkyl}$, $(\text{C}_2\text{--C}_8)\text{alkenyl}$, $(\text{C}_2\text{--C}_8)\text{alkynyl}$, benzyl, aryl, or $(\text{C}_2\text{--C}_5)\text{heteroaryl}$;

each occurrence of R^6 is independently H, $(\text{C}_1\text{--C}_8)\text{alkyl}$, $(\text{C}_2\text{--C}_8)\text{alkenyl}$, $(\text{C}_2\text{--C}_8)\text{alkynyl}$, benzyl, aryl, $(\text{C}_2\text{--C}_5)\text{heteroaryl}$, or $(\text{C}_0\text{--C}_8)\text{alkyl--C}(\text{O})\text{O--R}^5$ or the R^6 groups can join to form a heterocycloalkyl group; and

the * represents a chiral-carbon center.

11. (Original) A compound of claim 10, wherein R^1 is H, (C_1-C_4) alkyl, CH_2OCH_3 , $CH_2CH_2OCH_3$, or

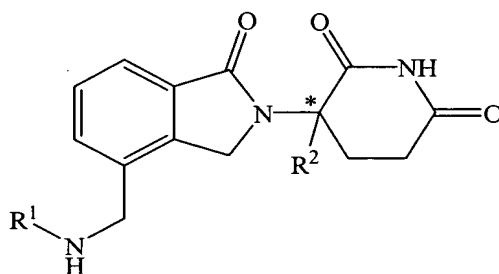


wherein Q is O or S, and each occurrence of R^7 is independently H, (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, halogen, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, (C_0-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, (C_1-C_8) alkyl- $O(CO)R^5$, or $C(O)OR^5$, or adjacent occurrences of R^7 can be taken together to form a bicyclic alkyl or aryl ring.

12. (Original) A compound of claim 10, wherein R^1 is $C(O)R^3$.

13. (Original) A compound of claim 10, wherein R^1 is $C(O)OR^4$.

14. (Previously presented) A compound having the formula:



wherein:

R^1 is H, (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, $C(O)R^3$, $C(S)R^3$, $C(O)OR^4$, (C_1-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, $C(O)NHR^3$, $C(S)NHR^3$, $C(O)NR^3R^3$, $C(S)NR^3R^3$ or (C_1-C_8) alkyl- $O(CO)R^5$;

R^2 is H or (C_1-C_8) alkyl;

R^3 and R^3' are independently (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, (C_0-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, (C_1-C_8) alkyl- $O(CO)R^5$, or $C(O)OR^5$;

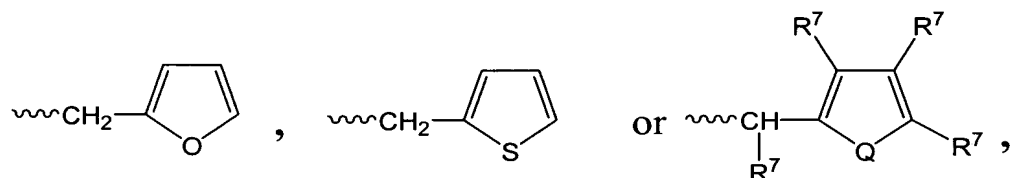
R^4 is (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, (C_1-C_4) alkyl- OR^5 , benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, or (C_0-C_4) alkyl- (C_2-C_5) heteroaryl;

R^5 is (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, or (C_2-C_5) heteroaryl;

each occurrence of R^6 is independently H, (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_2-C_5) heteroaryl, or (C_0-C_8) alkyl- $C(O)O-R^5$ or the R^6 groups can join to form a heterocycloalkyl group; and

the * represents a chiral-carbon center.

15. (Original) A compound of claim 14, wherein R^1 is H, (C_1-C_4) alkyl, CH_2OCH_3 , $CH_2CH_2OCH_3$, or

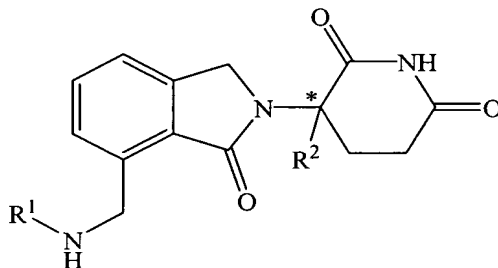


wherein Q is O or S, and each occurrence of R^7 is independently H, (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, halogen, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, (C_0-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, (C_1-C_8) alkyl- $O(CO)R^5$, or $C(O)OR^5$, or adjacent occurrences of R^7 can be taken together to form a bicyclic alkyl or aryl ring.

16. (Original) A compound of claim 14, wherein R^1 is $C(O)R^3$.

17. (Original) A compound of claim 14, wherein R^1 is $C(O)OR^4$.

18. (Previously presented) A compound having the formula:



wherein:

R^1 is H, (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, $C(O)R^3$, $C(S)R^3$, $C(O)OR^4$, (C_1-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, $C(O)NHR^3$, $C(S)NHR^3$, $C(O)NR^3R^{3'}$, $C(S)NR^3R^{3'}$ or (C_1-C_8) alkyl- $O(CO)R^5$;

R^2 is H or (C_1-C_8) alkyl;

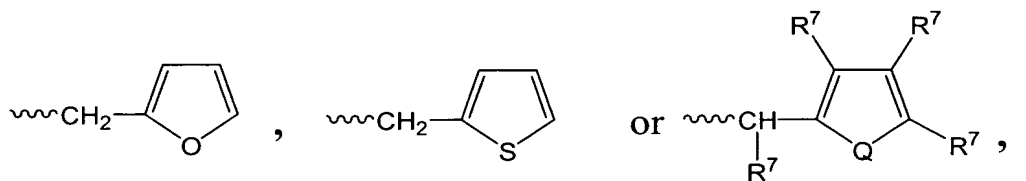
R^3 and $R^{3'}$ are independently (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, (C_0-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, (C_1-C_8) alkyl- $O(CO)R^5$, or $C(O)OR^5$;

R^4 is (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, (C_1-C_4) alkyl- OR^5 , benzyl, aryl, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, or (C_0-C_4) alkyl- (C_2-C_5) heteroaryl;

R^5 is (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, or (C_2-C_5) heteroaryl;

each occurrence of R^6 is independently H, (C_1-C_8) alkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, (C_2-C_5) heteroaryl, or (C_0-C_8) alkyl- $C(O)OR^5$ or the R^6 groups can join to form a heterocycloalkyl group; and
the * represents a chiral-carbon center.

19. (Original) A compound of claim 18, wherein R^1 is H, (C_1-C_4) alkyl, CH_2OCH_3 , $CH_2CH_2OCH_3$ or



wherein Q is O or S, and each occurrence of R^7 is independently H, (C_1-C_8) alkyl, (C_3-C_7) cycloalkyl, (C_2-C_8) alkenyl, (C_2-C_8) alkynyl, benzyl, aryl, halogen, (C_0-C_4) alkyl- (C_1-C_6) heterocycloalkyl, (C_0-C_4) alkyl- (C_2-C_5) heteroaryl, (C_0-C_8) alkyl- $N(R^6)_2$, (C_1-C_8) alkyl- OR^5 , (C_1-C_8) alkyl- $C(O)OR^5$, (C_1-C_8) alkyl- $O(CO)R^5$, or $C(O)OR^5$, or adjacent occurrences of R^7 can be taken together to form a bicyclic alkyl or aryl ring.

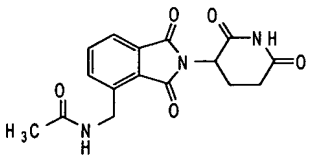
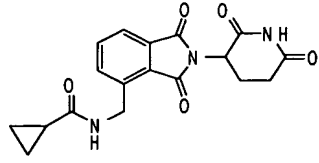
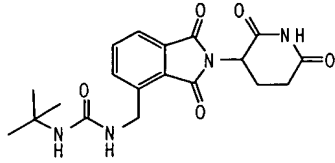
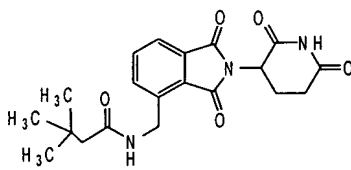
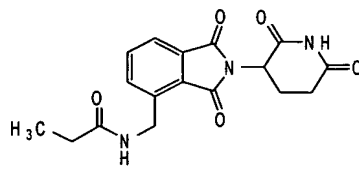
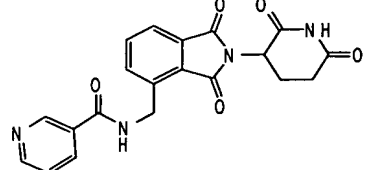
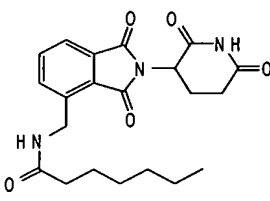
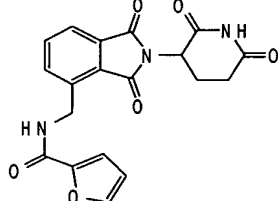
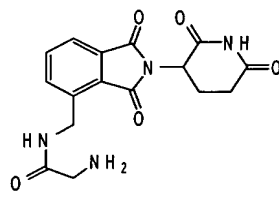
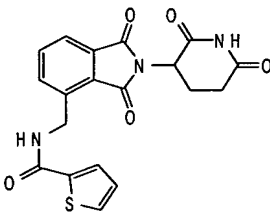
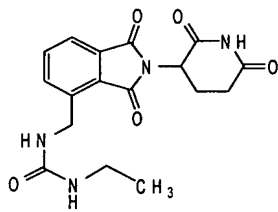
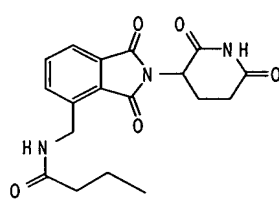
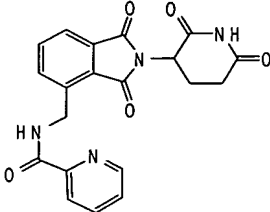
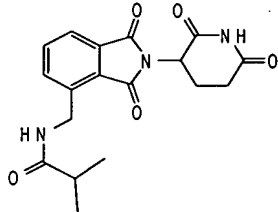
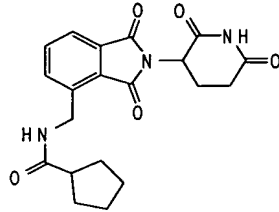
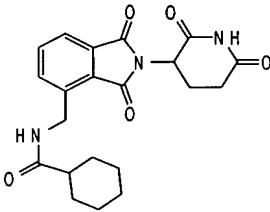
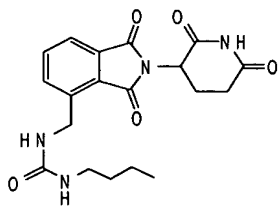
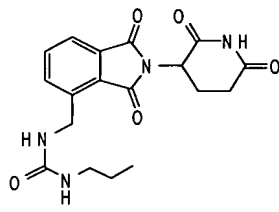
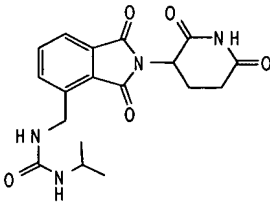
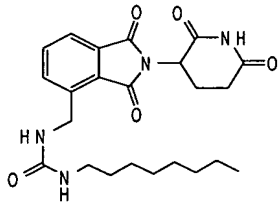
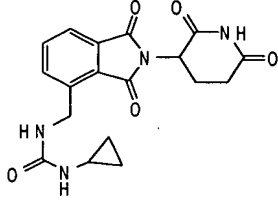
20. (Original) A compound of claim 18, wherein R^1 is $C(O)R^3$.

21. (Original) A compound of claim 18, wherein R¹ is C(O)OR⁴.

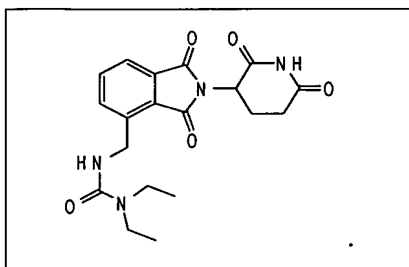
22-100. (Canceled).

101. (Currently amended) A compound of claim 10, which is: N-[2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-acetamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]cyclopropyl-carboxamide; 1-tert-butyl-3-[2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-urea; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-3,3-dimethylbutanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-propanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-3-pyridylcarboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]heptanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-2-furylcarboxamide; 2-amino-N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-acetamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-2-thienylcarboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(ethylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]butanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-2-pyridylcarboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]undecanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-2-methylpropanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]cyclopentylcarboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]cyclohexylcarboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(butylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(propylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(methylethylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(octylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(cyclopropylamino)carboxamide; or N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-(diethylamino)carboxamide.

(Claim 101 continued on next page)

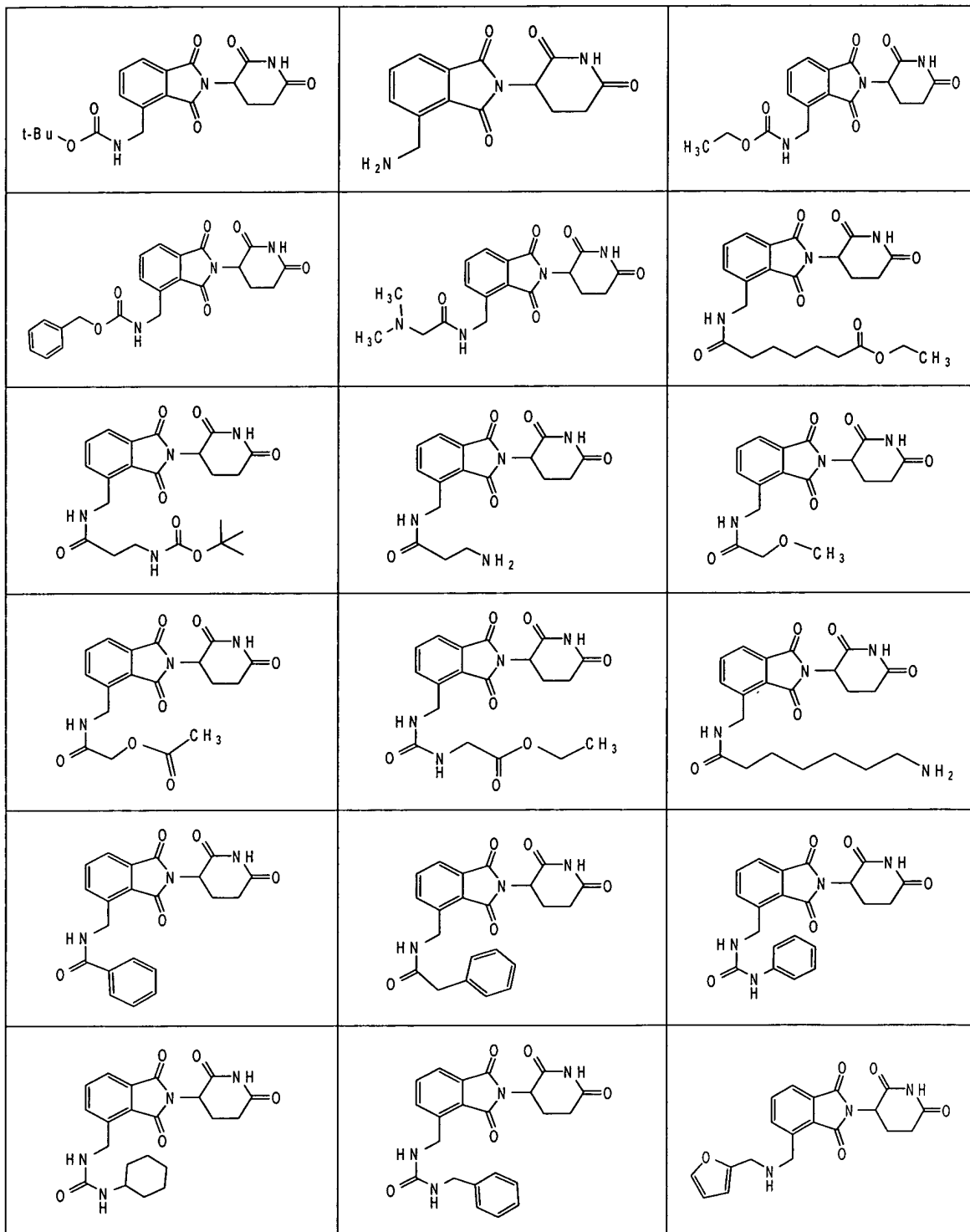
		
		
		
		
		
		
		

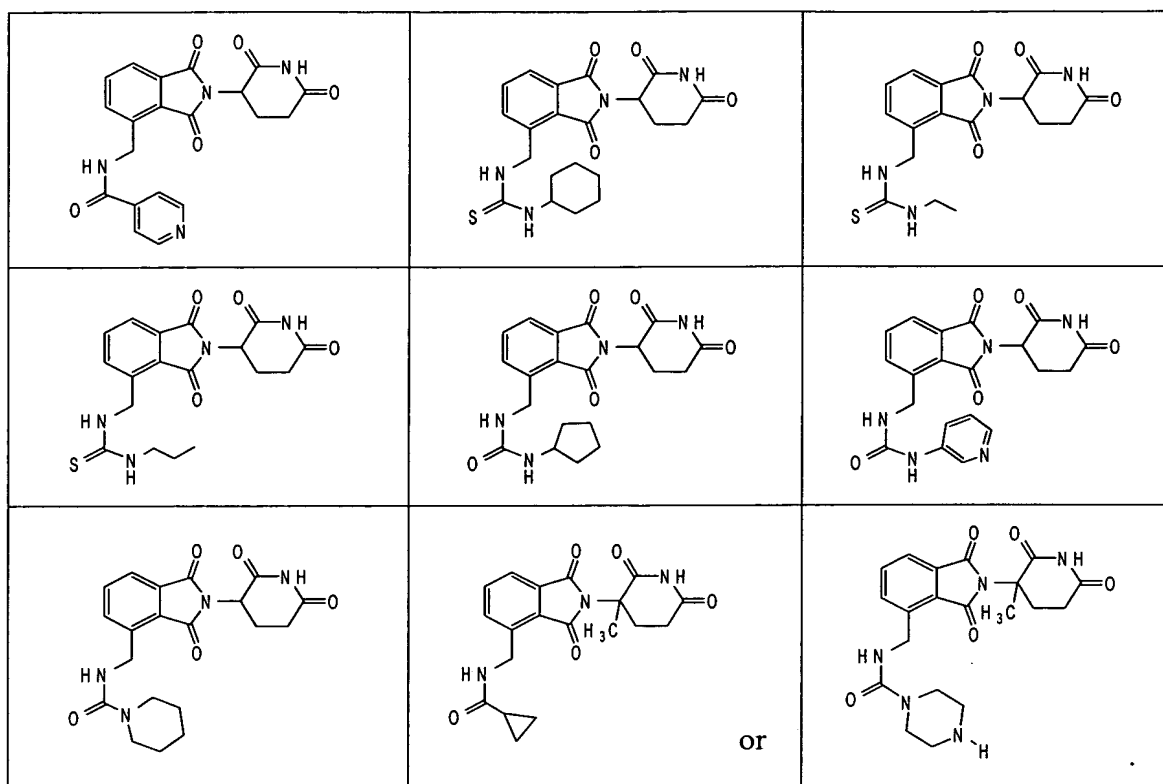
or



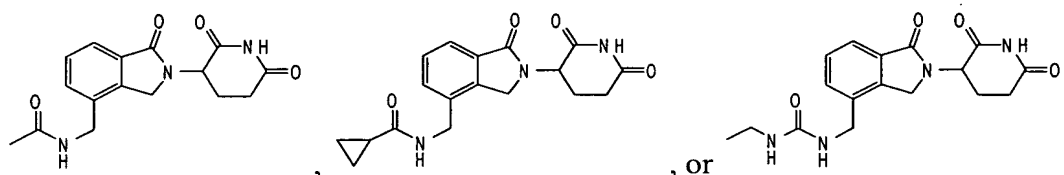
102. (Currently amended) A compound of claim 10, which is: ~~[2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-carbamic acid tert-butyl ester; 4-(aminomethyl)-2-(2,6-dioxo(3-piperidyl))-isoindoline-1,3-dione; [2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-carbamic acid ethyl ester; [2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-carbamic acid benzyl ester; 2-(dimethylamino)-N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]acetamide; ethyl 6-(3N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]carbamoyl)hexanoate; 3-[(tert-butoxy)carbonylamino]-N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]propanamide; 3-amino-N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]propanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]-2-methoxyacetamide; (N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]carbamoyl)methyl acetate; ethyl 2-[N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]carbamoyl]amino]acetate; 7-amino-N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]heptanamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl]benzamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl](phenylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl](benzylamino)carboxamide; 2-(2,6-dioxo-piperidin-3-yl)-4-[[[(furan-2-ylmethyl)-amino-methyl]-isoindole-1,3-dione]; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxo-2,3-dihydro-1H-isoindol-4-ylmethyl]-isonicotinamide; 2-(2,6-dioxo(3-piperidyl))-4-[[[(cyclohexylamino)thioxomethyl]amino]methyl]isoindole-1,3-dione; 2-(2,6-dioxo(3-piperidyl))-4-[[[(ethylamino)thioxomethyl]amino]methyl]isoindole-1,3-dione; 2-(2,6-dioxo(3-piperidyl))-4-[[[(propylamino)thioxomethyl]amino]methyl]isoindole-1,3-dione; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl](cyclopentylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-~~

dioxoisindolin-4-yl)methyl}(3-pyridylamino)carboxamide; N-[[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisindolin-4-yl)methyl]piperidylcarboxamide; or piperazine-1-carboxylic acid [2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isindol-4-ylmethyl]-amide.





103. (Currently amended) A compound of claim 14, which is: ~~N-[2-(2,6-dioxo-piperidin-3-yl)-1-oxo-2,3-dihydro-1H-isoindol-4-ylmethyl]acetamide; N-[[2-(2,6-dioxo(3-piperidyl))-1-oxoisoindolin-4-yl]methyl]cyclopropylcarboxamide; or N-[[2-(2,6-dioxo(3-piperidyl))-1-oxoisoindolin-4-yl]methyl](ethylamino)carboxamide.~~



104-105. (Canceled)